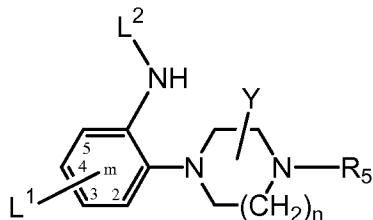


LISTING OF CLAIMS

1. (Original) A compound of formula (I):



formula (I)

and enantiomers, diastereomers and pharmaceutically acceptable salts thereof, wherein:
L¹ is a substituent moiety having a variable position “m”, wherein “m” represents a carbon atom number corresponding to a point of attachment for the L¹ substituent moiety on the anilino ring of formula (I);

L¹ is selected from the group consisting of R_{1b}, R₂-C(O), R_{1a}-SO₂ and R_{1a}-O(O)C-;

R_{1a} is C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

R_{1b} is C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

R₂ is heterocyclyl optionally substituted on a nitrogen atom with C₁₋₈alkyl;

L² is selected from the group consisting of R₃-C(O)-, R₄-SO₂-, R₆-NHC(S)- and R₆-NHC(O)-;

R₃ is selected from the group consisting of

(a) C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; wherein said aryl is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

- (b) aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- (c) heteroaryl optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro and aryl, wherein said aryl is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₄ is selected from the group consisting of

- (d) C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,
- (e) aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₆ is aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

- (f) C₁₋₈alkyl optionally substituted with one or more aryl substituents, wherein said aryl is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl; wherein said aryl' is optionally substituted with one or more substituents

independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

- (g) C₃₋₈cycloalkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino,

- mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy and nitro; and,
- (h) aryl optionally substituted with one or more substituents independently selected from the group consisting of C_{1-8} alkyl, C_{1-8} alkoxy, amino, mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy and nitro;

Y is one or more optionally present C_{1-8} alkyl substituents optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy, nitro, C_{3-8} cycloalkyl, aryl and heteroaryl, wherein said C_{3-8} cycloalkyl, aryl and heteroaryl are optionally further substituted;

m is an integer from 2 to 5 which represents the carbon atom number corresponding to the point of attachment for the L^1 substituent moiety on the anilino ring of formula (I); and, n is an integer from 1 to 2.

2. (Original) The compound of claim 1, wherein when L^2 is $R_3-C(O)-$ and R_3 is selected from the group consisting of unsubstituted C_{1-8} alkyl, substituted aryl, unsubstituted aryl, substituted heteroaryl and unsubstituted heteroaryl, then L^1 is $R_2-C(O)$.
3. (Original) The compound of claim 1, wherein when L^2 is $R_3-C(O)-$ and R_3 is selected from the group consisting of unsubstituted C_{1-8} alkyl, substituted aryl, unsubstituted aryl, substituted heteroaryl and unsubstituted heteroaryl, then R_5 is C_{1-8} alkyl optionally substituted with one or more optionally substituted aryl substituents.
4. (Original) The compound of claim 1, wherein when L^2 is R_4-SO_2- and R_4 is unsubstituted C_{1-8} alkyl, then L^1 is $R_2-C(O)$, wherein R_2 is substituted or unsubstituted heterocyclyl.
5. (Original) The compound of claim 1, wherein when L^2 is R_4-SO_2- and R_4 is unsubstituted C_{1-8} alkyl, then R_5 is C_{1-8} alkyl optionally substituted with one or more optionally substituted aryl substituents.

6. (Original) The compound of claim 1, wherein when L¹ is selected from the group consisting of R_{1b} and R_{1a}-O(O)C-, then L² is R₆-NHC(O)-, wherein R₆ is substituted or unsubstituted aryl.
7. (Original) The compound of claim 1, wherein when L¹ is selected from the group consisting of R_{1b} and R_{1a}-O(O)C-, then R₅ is C₁₋₈alkyl optionally substituted with one or more optionally substituted aryl substituents.
8. (Original) The compound of claim 1, wherein R_{1a} is C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy; R_{1b} is C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy; R₃ is selected from the group consisting of
 - (a) C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and, wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
 - (b) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
 - (c) heteroaryl optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro and aryl, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₄ is selected from the group consisting of

- (d) C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,
- (e) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₆ is aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

- (f) C₁₋₈alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl; wherein said aryl' is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- (g) C₃₋₈cycloalkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- (h) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

Y is one or two optionally present C₁₋₈alkyl substituents optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro,

C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted;

9. (Original) The compound of claim 1, wherein when L² is R₃-C(O)- and R₃ is selected from the group consisting of unsubstituted C₁₋₈alkyl, substituted aryl, unsubstituted aryl, substituted heteroaryl and unsubstituted heteroaryl, then R₅ is C₁₋₈alkyl optionally substituted with one or two optionally substituted aryl substituents.
10. (Original) The compound of claim 1, wherein when L² is R₄-SO₂- and R₄ is unsubstituted C₁₋₈alkyl, then R₅ is C₁₋₈alkyl optionally substituted with one or two optionally substituted aryl substituents.
11. (Original) The compound of claim 1, wherein when L¹ is selected from the group consisting of R_{1b} and R_{1a}-O(O)C-, then R₅ is C₁₋₈alkyl optionally substituted with one or two optionally substituted aryl substituents.
12. (Original) The compound of claim 1, wherein R_{1a} is C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, halogen and hydroxy; R_{1b} is C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, halogen and hydroxy; R₂ is piperazinyl optionally substituted on a nitrogen atom with C₁₋₄alkyl; L² is selected from the group consisting of R₃-C(O)-, R₄-SO₂- and R₆-NHC(O)-; R₃ is selected from the group consisting of
 - (a) C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl;
 - (b) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, halogen and hydroxy; and,

(c) heteroaryl optionally substituted on a secondary amine atom with C₁₋₄alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro and aryl, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₄ is selected from the group consisting of

- (d) C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,
- (e) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₆ is aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

- (f) C₁₋₄alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl; wherein said aryl' is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₄alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

- (g) C₃₋₈cycloalkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

- (h) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

Y is absent;

m is an integer from 3 to 4 which represents the carbon atom number corresponding to the point of attachment for the L¹ substituent moiety on the anilino ring of formula (I); and, n is 1.

13. (Original) The compound of claim 12, wherein

R_{1a} is C₁₋₄alkyl;

R_{1b} is hydroxy(C₁₋₄)alkyl-;

R₃ is selected from the group consisting of

- (a) C₁₋₄alkyl;
(b) phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, halogen and hydroxy; and,
(c) furyl optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and aryl, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₄ is selected from the group consisting of

- (d) C₁₋₄alkyl; and,
(e) phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₆ is phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

R₅ is selected from the group consisting of

- (f) C₁₋₄alkyl optionally substituted with one or two aryl substituents, wherein aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl;
(g) C₃₋₈cycloalkyl; and,

(h) aryl.

14. (Original) The compound of claim 13, wherein R₃ is selected from the group consisting of

- (a) C₁₋₄alkyl;
- (b) phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl and halogen; and,
- (c) furyl optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl and phenyl; wherein said phenyl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl and halogen;

R₄ is selected from the group consisting of

- (d) C₁₋₄alkyl; and,
- (e) phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl and halogen;

R₆ is phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, amino, halogen and hydroxy; and,

R₅ is C₁₋₄alkyl optionally substituted with one or two phenyl substituents, wherein phenyl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, amino, halogen and hydroxy.

15. (Original) The compound of claim 1, wherein R_{1a} is C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy.

16. (Original) The compound of claim 1, wherein R_{1a} is C₁₋₄alkyl optionally substituted with one substituent selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy.

17. (Original) The compound of claim 1, wherein R_{1a} is C₁₋₄alkyl.

18. (Original) The compound of claim 1, wherein R_{1b} is C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino and hydroxy.

19. (Original) The compound of claim 1, wherein R_{1b} is C₁₋₄alkyl optionally

- substituted with one substituent selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy.
20. (Original) The compound of claim 1, wherein R_{1b} is C₁₋₄alkyl optionally substituted with hydroxy.
21. (Original) The compound of claim 1, wherein R₂ is piperazinyl optionally substituted on a nitrogen atom with C₁₋₄alkyl.
22. (Original) The compound of claim 1, wherein L² is R₃-C(O)-.
23. (Original) The compound of claim 22, wherein R₃ is selected from the group consisting of
- (a) C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl;
 - (b) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, halogen and hydroxy;
 - (c) heteroaryl optionally substituted on a secondary amine atom with C₁₋₄alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro and aryl, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro.
24. (Original) The compound of claim 22, wherein R₃ is selected from the group consisting of
- (a) C₁₋₄alkyl;
 - (b) phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, halogen and hydroxy; and,
 - (c) furyl optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and aryl, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro.

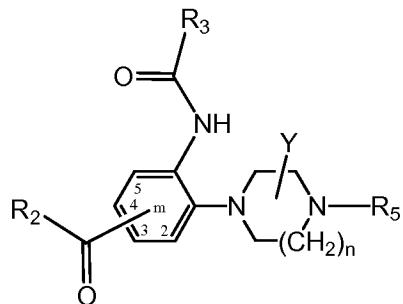
25. (Original) The compound of claim 22, wherein R₃ is selected from the group consisting of
 - (a) C₁₋₄alkyl;
 - (b) phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl and halogen; and,
 - (c) furyl optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl and phenyl; wherein said phenyl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl and halogen.
26. (Original) The compound of claim 1, wherein L² is R₃-C(O)- and R₅ is C₁₋₈alkyl optionally substituted with one or two optionally substituted aryl substituents.
27. (Original) The compound of claim 26, wherein R₃ is selected from the group consisting of
 - (a) C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl;
 - (b) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, halogen and hydroxy;
 - (c) heteroaryl optionally substituted on a secondary amine atom with C₁₋₄alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro and aryl, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro.
28. (Original) The compound of claim 1, wherein L² is R₄-SO₂-.
29. (Original) The compound of claim 28, wherein R₄ is selected from the group consisting of
 - (d) C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,

- (e) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro.
30. (Original) The compound of claim 28, wherein R₄ is selected from the group consisting of
- (d) C₁₋₄alkyl; and,
 - (e) phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl and halogen.
31. (Original) The compound of claim 1, wherein L² is R₄-SO₂- and R₅ is C₁₋₈alkyl optionally substituted with one or two optionally substituted aryl substituents.
32. (Original) The compound of claim 31, wherein R₄ is selected from the group consisting of
- (d) C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,
 - (e) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro.
33. (Original) The compound of claim 1, wherein L² is R₆-NHC(O)-.
34. (Original) The compound of claim 33, wherein R₆ is phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro.
35. (Original) The compound of claim 1, wherein L² is R₆-NHC(O)- and R₅ is C₁₋₈alkyl optionally substituted with one or two optionally substituted aryl substituents.
36. (Original) The compound of claim 35, wherein R₆ is phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro.
37. (Original) The compound of claim 1, wherein R₅ is selected from the group consisting of

- (f) C_{1-4} alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C_{1-4} alkyl, C_{1-4} alkoxy, amino, mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl;
 - (g) C_{3-8} cycloalkyl optionally substituted with one or two substituents independently selected from the group consisting of C_{1-4} alkyl, C_{1-4} alkoxy, amino, mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy and nitro; and,
 - (h) aryl optionally substituted with one or two substituents independently selected from the group consisting of C_{1-4} alkyl, C_{1-4} alkoxy, amino, mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy and nitro.
38. (Original) The compound of claim 1, wherein R_5 is selected from the group consisting of
- (f) C_{1-4} alkyl optionally substituted with one or two aryl substituents, wherein aryl is optionally substituted with one or two substituents independently selected from the group consisting of C_{1-4} alkyl, C_{1-4} alkoxy, amino, mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy and nitro;
 - (g) C_{3-8} cycloalkyl; and,
 - (h) aryl.

39. (Original) The compound of claim 1, wherein R_5 is C_{1-4} alkyl optionally substituted with one or two phenyl substituents, wherein phenyl is optionally substituted with one or two substituents independently selected from the group consisting of C_{1-4} alkyl, amino, halogen and hydroxy.

40. (Original) The compound of claim 1, wherein the compound of formula (I) is selected from a compound of formula (Ia):



formula (Ia)

and enantiomers, diastereomers and pharmaceutically acceptable salts thereof, wherein:

R₂-C(O)- is a substituent moiety having a variable position “m”, wherein “m” represents a carbon atom number corresponding to a point of attachment for the R₂-C(O)- substituent moiety on the anilino ring of formula (Ia);

R₂ is heterocyclyl optionally substituted on a nitrogen atom with C₁₋₈alkyl;

R₃ is selected from the group consisting of

- (a) C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl;
wherein said aryl is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- (b) aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- (c) heteroaryl optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro and aryl, wherein said aryl is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

- (f) C₁₋₈alkyl optionally substituted with one or more aryl substituents, wherein said aryl is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl;
wherein said aryl' is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

- wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- (g) C₃₋₈cycloalkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- (h) aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- Y is one or more optionally present C₁₋₈alkyl substituents optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted;
- m is an integer from 2 to 5 which represents the carbon atom number corresponding to the point of attachment for the R₂-C(O)- substituent moiety on the anilino ring of formula (Ia); and, n is an integer from 1 to 2.
41. (Original) The compound of claim 40, wherein R₃ is selected from the group consisting of
- (a) C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and, wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- (b) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- (c) heteroaryl optionally substituted on a secondary amine atom with C₁₋₈alkyl, and

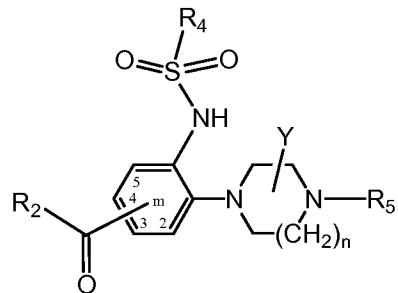
optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro and aryl, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

- (f) C₁₋₈alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl; wherein said aryl' is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- (g) C₃₋₈cycloalkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- (h) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

Y is one or two optionally present C₁₋₈alkyl substituents optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted.

42. (Original) The compound of claim 1, wherein the compound of formula (I) is selected from a compound of formula (Ib):



formula (Ib)

and enantiomers, diastereomers and pharmaceutically acceptable salts thereof, wherein: R₂-C(O)- is a substituent moiety having a variable position "m", wherein "m" represents a carbon atom number corresponding to a point of attachment for the R₂-C(O)- substituent moiety on the anilino ring of formula (Ib);

R₂ is heterocyclyl optionally substituted on a nitrogen atom with C₁₋₈alkyl;

R₄ is selected from the group consisting of

- (d) C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,
(e) aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

- (f) C₁₋₈alkyl optionally substituted with one or more aryl substituents, wherein said aryl is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl; wherein said aryl' is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of

C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

- (g) C₃₋₈cycloalkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- (h) aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

Y is one or more optionally present C₁₋₈alkyl substituents optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted;

m is an integer from 2 to 5 which represents the carbon atom number corresponding to the point of attachment for the R₂-C(O)- substituent moiety on the anilino ring of formula (Ib); and, n is an integer from 1 to 2.

43. (Original) The compound of claim 42, wherein

R₄ is selected from the group consisting of

- (d) C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,
- (e) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

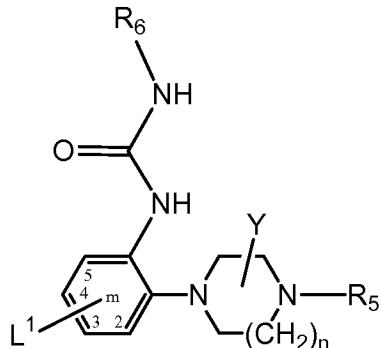
- (f) C₁₋₈alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl; wherein said aryl' is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

- (g) C₃₋₈cycloalkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- (h) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

Y is one or two optionally present C₁₋₈alkyl substituents optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted.

44. (Original) The compound of claim 1, wherein the compound of formula (I) is selected from a compound of formula (Ic):



formula (Ic)

and enantiomers, diastereomers and pharmaceutically acceptable salts thereof, wherein:
L¹ is a substituent moiety having a variable position "m", wherein "m" represents a carbon atom number corresponding to a point of attachment for the L¹ substituent moiety on the anilino ring of formula (Ic);

L¹ is selected from the group consisting of R_{1b}, R_{1a}-SO₂- and R_{1a}-O(O)C-;

R_{1a} is C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

R_{1b} is C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

R₆ is aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

- (f) C₁₋₈alkyl optionally substituted with one or more aryl substituents, wherein said aryl is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl; wherein said aryl' is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

- (g) C₃₋₈cycloalkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

- (h) aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

Y is one or more optionally present C₁₋₈alkyl substituents optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and

heteroaryl are optionally further substituted;

m is an integer from 2 to 5 which represents the carbon atom number corresponding to the point of attachment for the L¹ substituent moiety on the anilino ring of formula (Ic); and, n is an integer from 1 to 2.

45. (Original) The compound of claim 44, wherein

R_{1a} is C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

R_{1b} is C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

R₆ is aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

(f) C₁₋₈alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl; wherein said aryl' is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

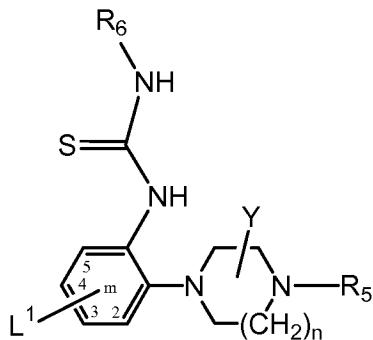
wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

(g) C₃₋₈cycloalkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

(h) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

Y is one or two optionally present C₁₋₈alkyl substituents optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted.

46. (Original) The compound of claim 1, wherein the compound of formula (I) is selected from a compound of formula (Id):



formula (Id)

and enantiomers, diastereomers and pharmaceutically acceptable salts thereof, wherein:
L¹ is a substituent moiety having a variable position "m", wherein "m" represents a carbon atom number corresponding to a point of attachment for the L¹ substituent moiety on the anilino ring of formula (Id);

L¹ is selected from the group consisting of R_{1b}, R₂-C(O)-, R_{1a}-SO₂- and R_{1a}-O(O)C-;

R_{1a} is C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

R_{1b} is C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

R₂ is heterocyclyl optionally substituted on a nitrogen atom with C₁₋₈alkyl;

R₆ is aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

- (f) C₁₋₈alkyl optionally substituted with one or more aryl substituents, wherein said aryl is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl; wherein said aryl' is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and, wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- (g) C₃₋₈cycloalkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- (h) aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

Y is one or more optionally present C₁₋₈alkyl substituents optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted;

m is an integer from 2 to 5 which represents the carbon atom number corresponding to the point of attachment for the L¹ substituent moiety on the anilino ring of formula (Id); and, n is an integer from 1 to 2.

47. (Original) The compound of claim 46, wherein R_{1a} is C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

R_{1b} is C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

R₆ is aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

(f) C₁₋₈alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl; wherein said aryl' is optionally substituted with one or two substituents

independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

(g) C₃₋₈cycloalkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

(h) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

Y is one or two optionally present C₁₋₈alkyl substituents optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted.

48. (Original) A compound selected from the group consisting of:

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-4-methyl-benzamide;

5-(4-chlorophenyl)-N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-2-methyl-3-furancarboxamide;

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-2-furancarboxamide;

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-propanamide;

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-4-methylbenzenesulfonamide;

4-chloro-*N*-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-benzenesulfonamide;

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-1-butanesulfonamide;

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-methanesulfonamide;

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(methylsulfonyl)phenyl]-*N'*-phenyl-urea.

N-[2-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-5-(hydroxymethyl)phenyl]-*N'*-phenyl-urea; and,

4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[(phenylamino)carbonyl]amino]-benzoic acid methyl ester.

49. (Original) A composition comprising a pharmaceutically acceptable carrier, excipient, tableting ingredient or diluent and the compound of claim 1.

50-55. Canceled.